

# *STATISTICAL PHYSICS FOR ADAPTIVE DISTRIBUTED CONTROL*

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## THE GOLDEN RULE

DO NOT:

*Find a value of a variable  $x$ ,  
that optimizes a function*

INSTEAD:

*Find a distribution over  $x$ ,  
that optimizes an expectation value*



## ADVANTAGES

- 1) *Arbitrary data types.*
  - 2) *Leverages continuous-space optimization.*  
*(“Gradient descent for symbolic variables”.)*
  - 3) *Akin to interior point methods.*
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- *Deep connections with statistical physics and game theory. So*
    - *Especially suited for distributed domains.*
    - *Especially suited for very large problems.*



# ROADMAP

1) *What is distributed control, formally?*



2) *Review information theory*



3) *Optimal control policy for distributed agents*



4) *How to find that policy in a distributed way*



## WHAT IS DISTRIBUTED CONTROL?

- 1) A set of  $N$  agents: Joint move  $x = (x_1, x_2, \dots, x_N)$
- 2) Since they are distributed, their joint probability is a product distribution:

$$q(x) = \prod_i q_i(x_i)$$

- This definition of distributed agents is adopted from (extensive form) noncooperative game theory.



## **EXAMPLE: KSAT**

- $x = \{0, 1\}^N$
- A set of many disjunctions, “clauses”, each involving  $K$  bits.  
E.g.,  $(x_2 \vee x_6 \vee \neg x_7)$  is a clause for  $K = 3$
- Goal: Find a bit-string  $x$  that simultaneously satisfies all clauses.  $G(x)$  is #violated clauses.
- For us, this goal becomes: find a  $q(x) = \prod_i q_i(x_i)$  tightly centered about such an  $x$ .

*The canonical computationally difficult problem*



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## ***REVIEW OF INFORMATION THEORY***

- 1) Want a quantification of how “uncertain” you are that you will observe a value  $i$  generated from  $P(i)$ .**
- 2) Require the uncertainty at seeing the IID pair  $(i, i')$  to equal the sum of the uncertainties for  $i$  and for  $i'$**
- 3) This forces the definition**

$$\text{uncertainty}(i) = -\ln[P(i)]$$



## REVIEW OF INFORMATION THEORY - 2

4) So expected uncertainty is the *Shannon entropy*

$$S(P) \propto -\sum_i P(i) \ln[P(i)]$$

- Concave over P
- $\alpha(P)$  is infinite at border of space of all P

5) *Information* in P,  $I(P)$ , is what's left after the uncertainty is removed:  $-S(P)$ .



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## ITERATIVE DISTRIBUTED CONTROL

- 1)  $s$  is current uncertainty of what  $x$  to pick, i.e., uncertainty of where  $q(x)$  is concentrated.
  - Early in the control process, high uncertainty.
- 2) Find  $q$  minimizing  $E_q(G)$  while consistent with  $s$ .
- 3) Reduce  $s$ . Return to (2).
- 4) Terminate at a  $q$  with good (low)  $E_q(G)$ .

*Can do (2) $\alpha$  (3) without ever explicitly specifying  $s$*



## ITERATIVE DISTRIBUTED CONTROL - 2

- 1) The central step is to “find the  $q$  that has lowest  $E_q(G)$  while consistent with  $S(q) = s$ ”.
- 2) So we must find the critical point of the Lagrangian
$$L(q, T) = E_q(G) + T[s - S(q)] ,$$
i.e., find the  $q$  and  $T$  such that  $\alpha L / \alpha q = \alpha L / \alpha T = 0$ 
  - Deep connections with statistical physics ( $L$  is “free energy” in mean-field theory), economics
- 3) Then we reduce  $s$ ; repeat (find next critical point).



## EXAMPLE: KSAT

$$1) S(q) = -\sum_i [b_i \ln(b_i) + (1 - b_i) \ln(1 - b_i)]$$

where  $b_i$  is  $q_i(x_i = \text{TRUE})$

$$\begin{aligned} 2) E_q(G) &= \sum_{\text{clauses } j, x} q(x) K_j(x) \\ &= \sum_{\text{clauses } j, x, i} \prod_i q_i(x_i) K_j(x) \end{aligned}$$

where  $K_j(x) = 1$  iff  $x$  violates clause  $j$

**Our algorithm:** i) Find  $q$  minimizing  $E_q(G) - \text{TS}(q)$ ;  
ii) Lower  $T$  and return to (i).



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## DISTRIBUTED SEARCH FOR $q$

So control reduces to finding  $q$  such that  $\alpha L / \alpha q = 0$

- 1) Since the agents make their moves in a distributed way, that  $q$  is a product distribution.
- 2) But they must also find that  $q$  in a distributed way.
- 3) There are two cases to consider:
  - i) Know functional form of  $G$ .
  - ii) Don't know functional form of  $G$  - must sample.



## MINIMIZING $L(q)$ VIA GRADIENT DESCENT

- 1) Each  $i$  works to minimize  $L(q_i, q_{(i)})$  using only partial information of the other agents' distribution,  $q_{(i)}$ .
- 2) The  $q_i(x_i)$  component of  $\alpha L(q)$ , projected onto the space of allowed  $q_i(x_i)$ , is

$$\frac{\alpha E_{q_{(i)}}(G | x_i) + \ln(q_i(x_i))}{\alpha x \alpha [\alpha E_{q_{(i)}}(G | x_i) + \ln(q_i(x_i))]}$$

- The subtracted term ensures  $q$  stays normalized



## GRADIENT DESCENT - 2

- 3) Each agent  $i$  knows its value of  $\ln(q_i(x_i))$ .
- 4) Each agent  $i$  knows the  $E_{q(i)}(G | x_i)$  terms.

Each agent knows how it should change its  $q_i$  under gradient descent over  $L(q)$

- 5) Gradient descent, even for categorical variables (!), and done in a distributed way.
- 6) Similarly the Hessian can readily be estimated (for Newton's method), etc.



## **EXAMPLE: KSAT**

1) Evaluate  $E_{q(i)}(G \mid x_i)$  - the expected number of violated clauses if bit  $i$  is in state  $x_i$  - for every  $i, x_i$

2) In gradient descent, decrease each  $q_i(x_i)$  by

$$\alpha [E_{q(i)}(G \mid x_i) + T \ln[q_i(x_i)] - \text{const}_j]$$

where  $\alpha$  is the stepsize, and  $\text{const}_j$  is an easy-to-evaluate normalization constant.

3) We actually have a different  $T$  for each clause, and adaptively update all of them.



## ADAPTIVE DISTRIBUTED CONTROL

1) In *adaptive* control, don't know functional form of  $G(x)$ . So use Monte Carlo:

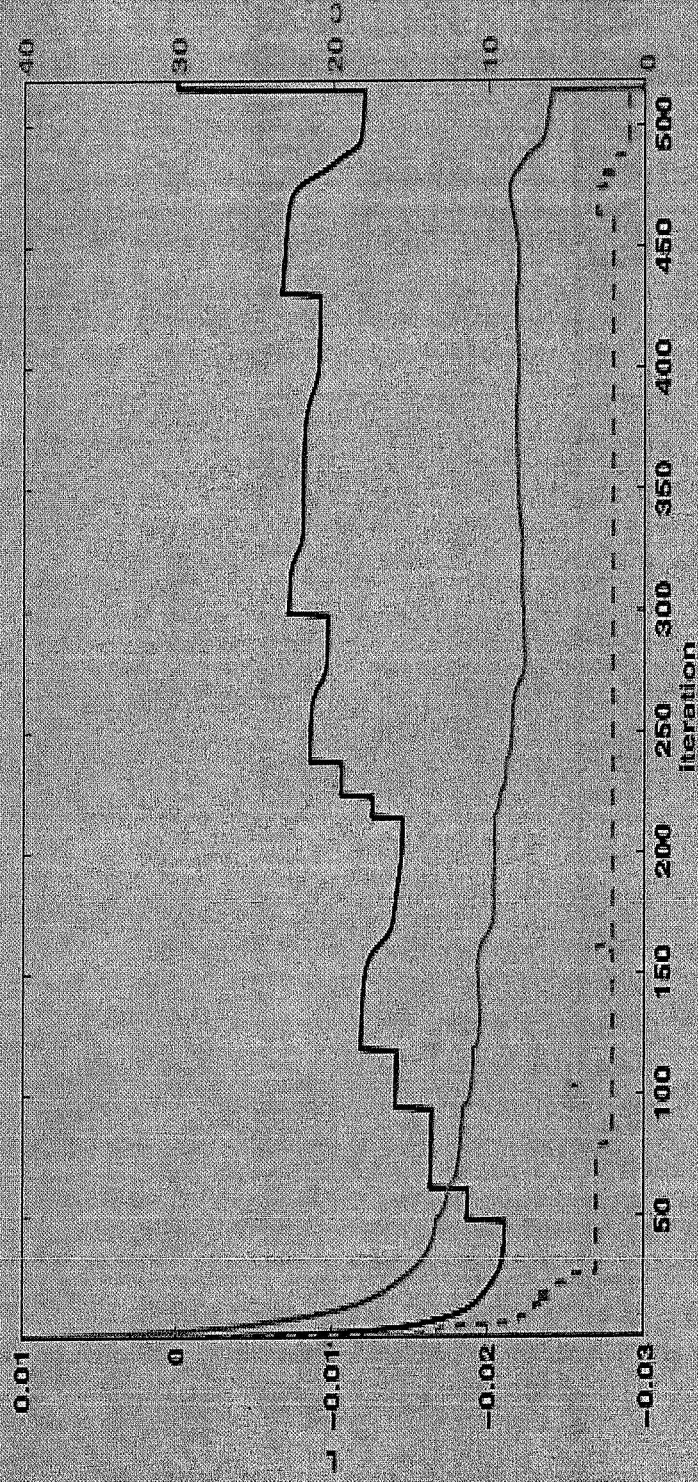
- Sample  $G(x)$  repeatedly according to  $q_i$ ;
- Each  $i$  independently estimates  $E_{q(i)}(G \mid x_i)$  for all its moves  $x_i$ ;
- Only 1 MC process, no matter how many agents

So each  $q_i$  can adaptively estimate its update



## **EXAMPLE: KSAT**

- i) Top plot is Lagrangian value vs. iteration;
- ii) Middle plot is average (under  $q$ ) number of constraint violations;
- iii) Bottom plot is mode (under  $q$ ) number of constraint violations.





## CONCLUSION

- 1) A distributed system is governed by a product distribution  $q$ , by definition.*
- 2) So distributed adaptive control is adaptive search for the  $q$  that optimizes  $E_q(G)$ .*
- 3) That search can be done many ways, e.g., gradient descent, with or without Monte Carlo sampling.*